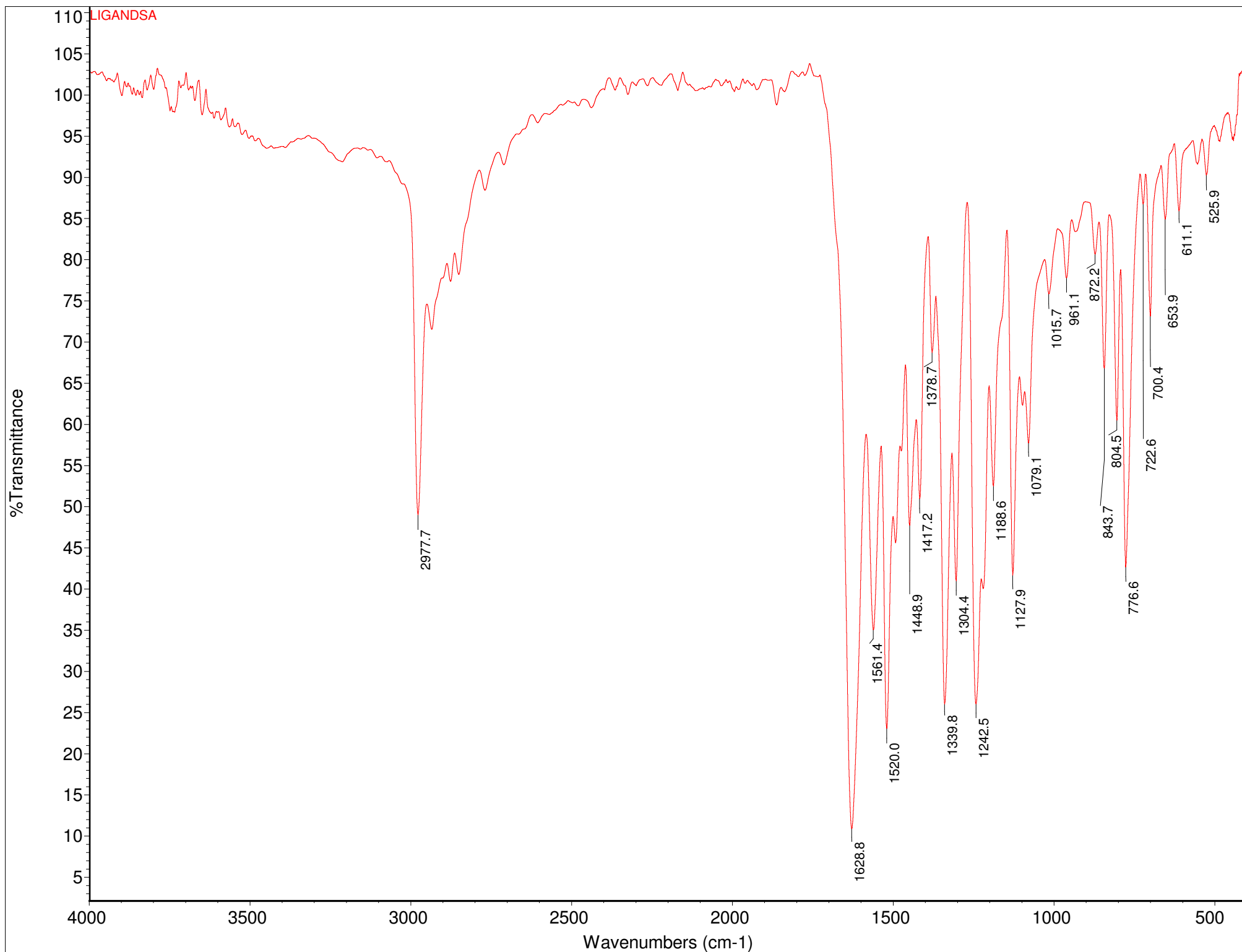
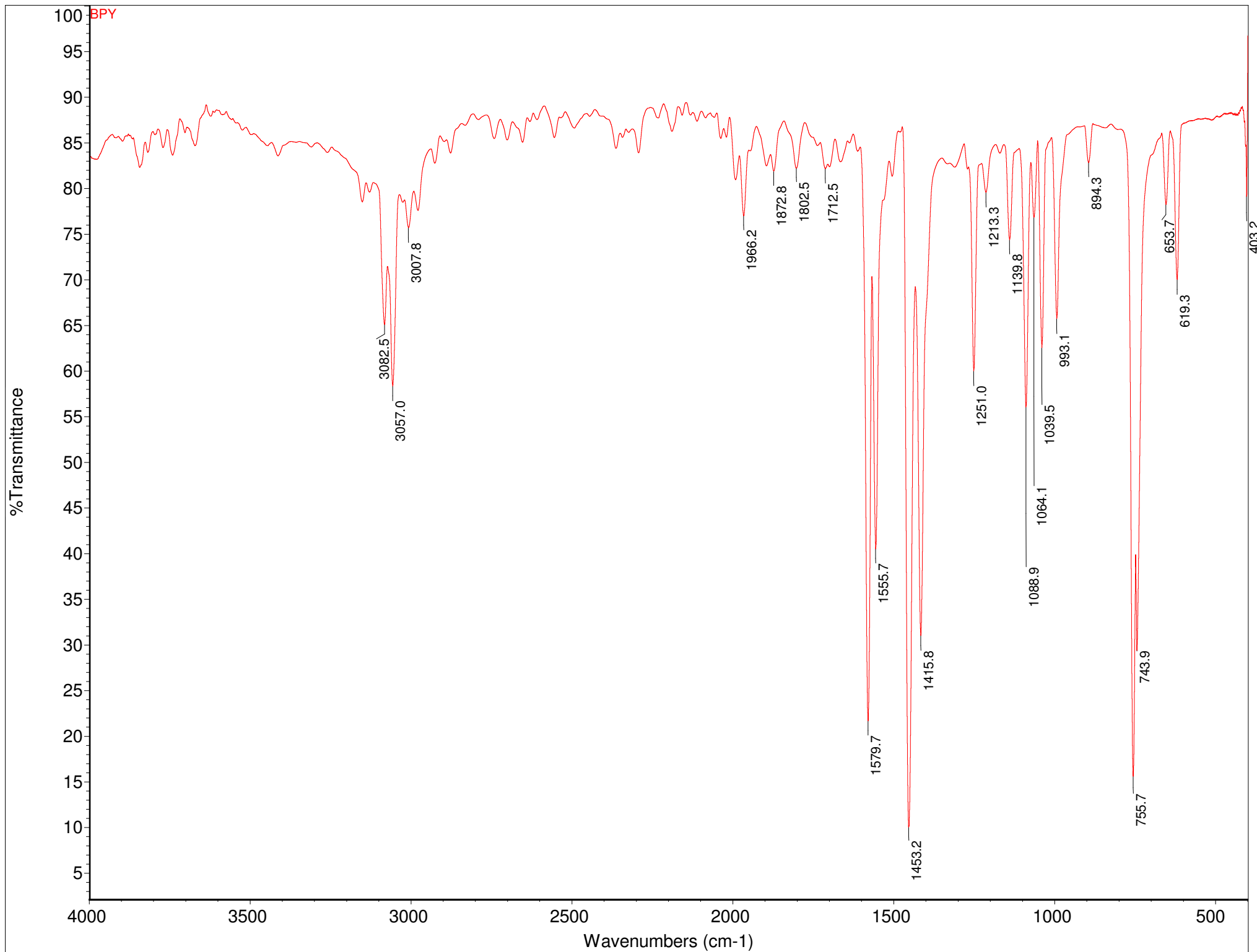
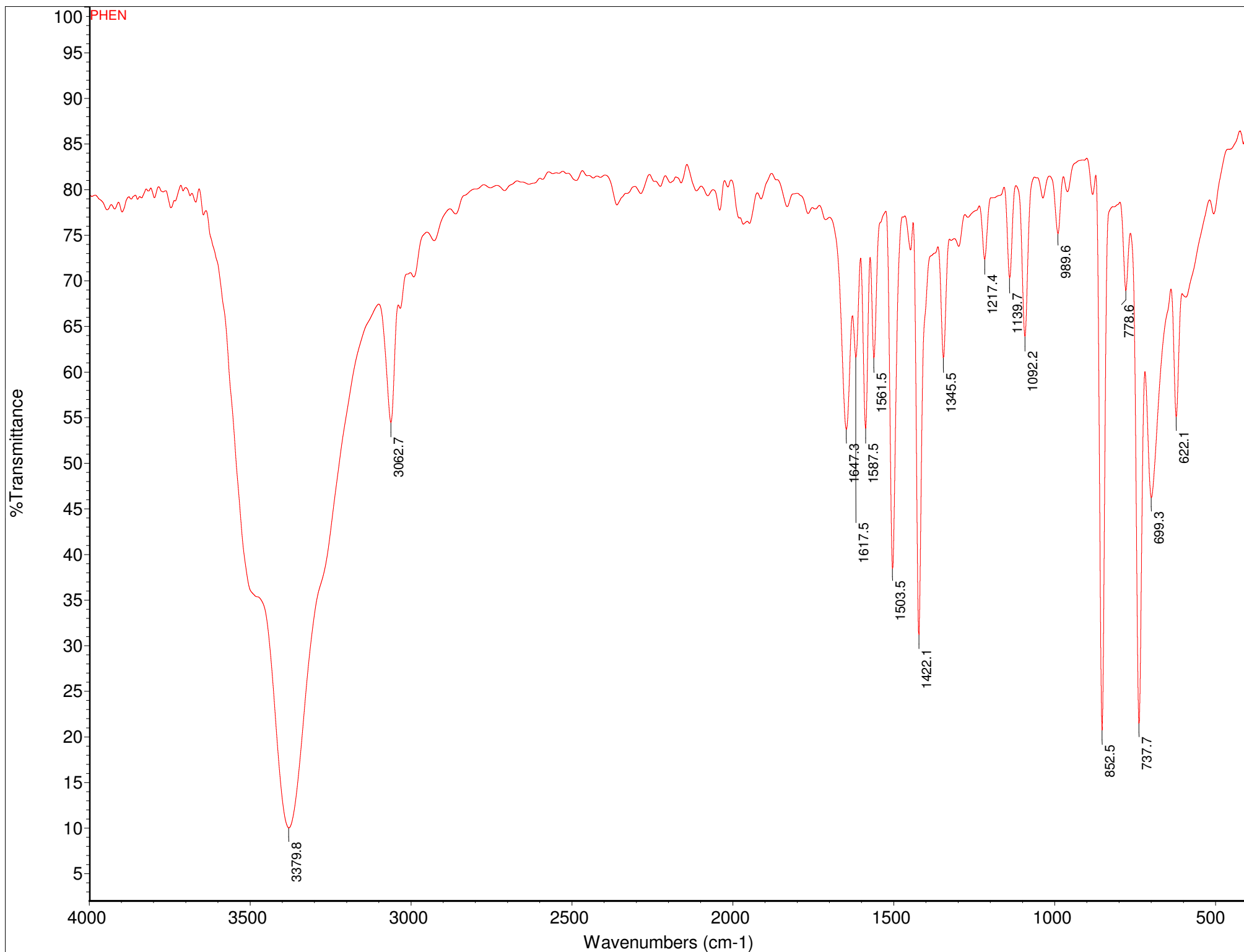


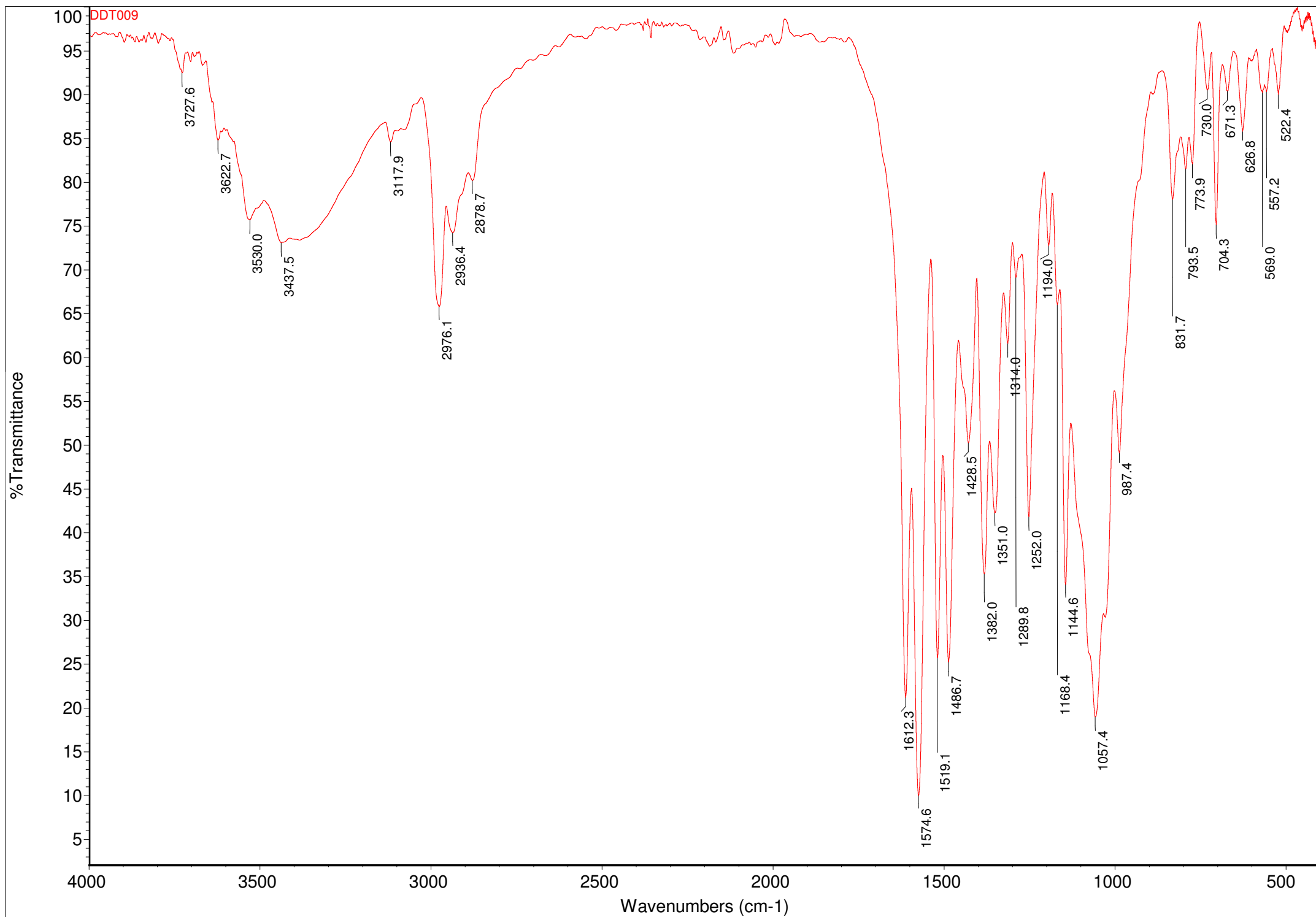
Supplementary data for the article:

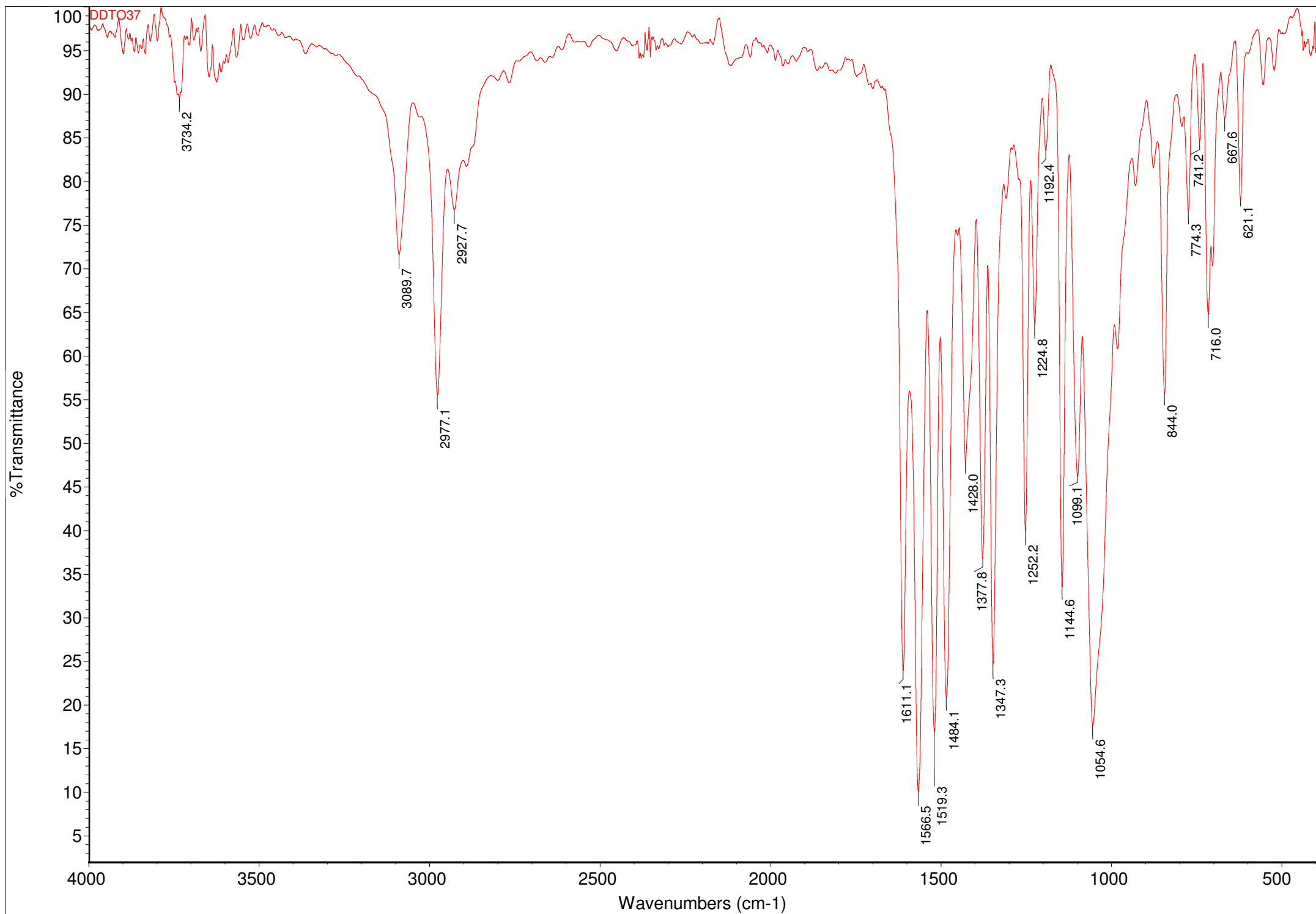
Dimitrijević, T.; Novaković, I.; Radanović, D.; Novaković, S. B.; Rodić, M. V.; Anđelković, K.; Šumar-Ristović, M. Synthesis, Spectral and Structural Characterization and Biological Activity of Cu(II) Complexes with 4-(Diethylamino)Salicylaldehyde and  $\alpha$ -Diimines. *Journal of Coordination Chemistry* **2020**, 73 (4), 702–716. <https://doi.org/10.1080/00958972.2020.1740212>











# Qualitative Compound Report

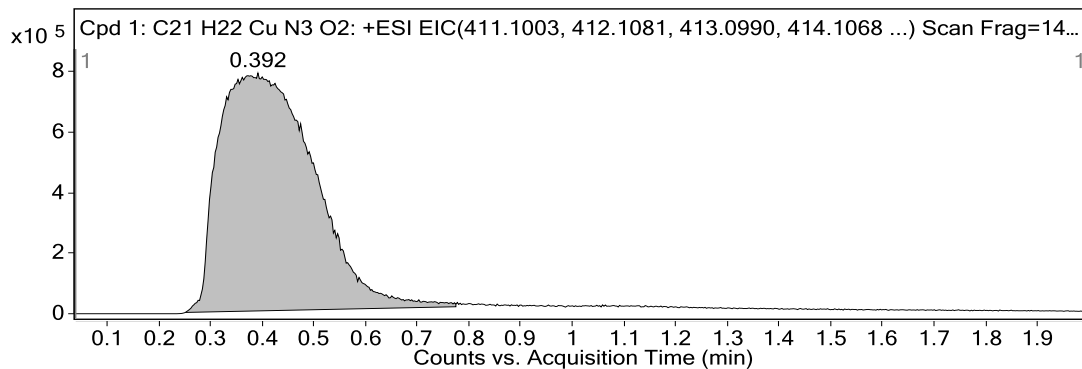
<b>Data File</b>	MJ_MS_1DMSO_MK_140V_pos1.d	<b>Sample Name</b>	1DMSO
<b>Sample Type</b>	Sample	<b>Position</b>	Vial 5
<b>Instrument Name</b>	DE1517B001	<b>User Name</b>	
<b>Acq Method</b>	Odredjivanje MM_MK_140V_pos.m	<b>Acquired Time</b>	11/29/2019 2:03:46 PM
<b>IRM Calibration Status</b>	Success	<b>DA Method</b>	EURL_target_PEST017_NNP_10V.m
<b>Comment</b>			

<b>Sample Group</b>		<b>Info.</b>	
<b>Stream Name</b>	LC 1	<b>Acquisition SW Version</b>	6200 series TOF/6500 series Q-TOF B.06.01 (B6157)

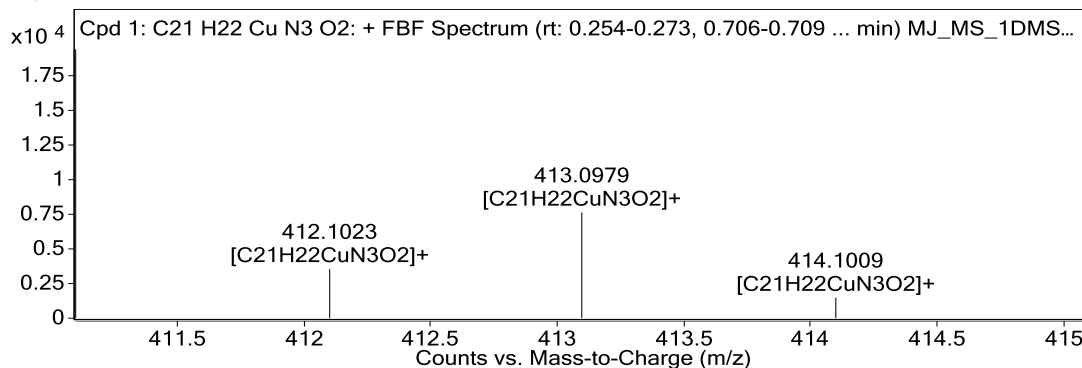
## Compound Table

Compound Label	RT	Mass	Abund	Formula	Tgt Mass	Diff (ppm)
Cpd 1: C21 H22 Cu N3 O2	0.392	411.0998	19387	C21 H22 Cu N3 O2	411.1008	-2.34

Compound Label	m/z	RT	Algorithm	Mass
Cpd 1: C21 H22 Cu N3 O2	411.0993	0.392	Find By Formula	411.0998

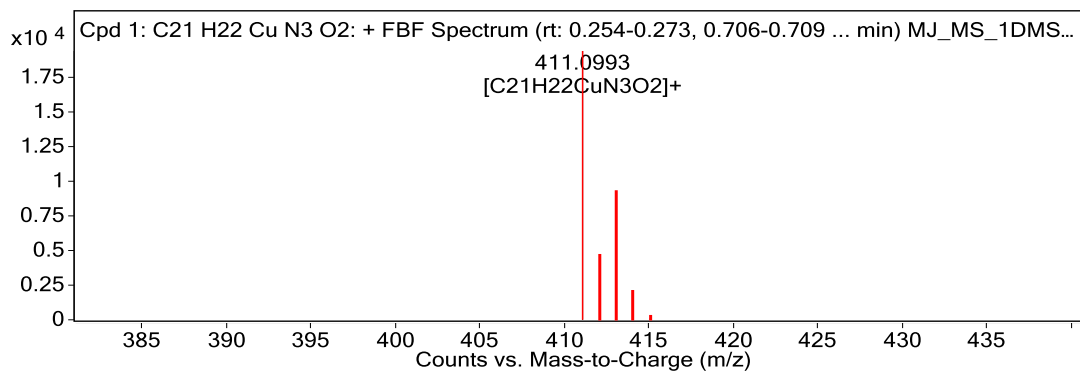


## MS Spectrum



## MS Zoomed Spectrum

# Qualitative Compound Report



## MS Spectrum Peak List

m/z	z	Abund	Formula	Ion
411.0993	1	19386.78	C <sub>21</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
412.1023	1	3551.74	C <sub>21</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
413.0979	1	7623.97	C <sub>21</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
414.1009	1	1477.12	C <sub>21</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
415.1028	1	115.38	C <sub>21</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>

--- End Of Report ---



# Qualitative Compound Report

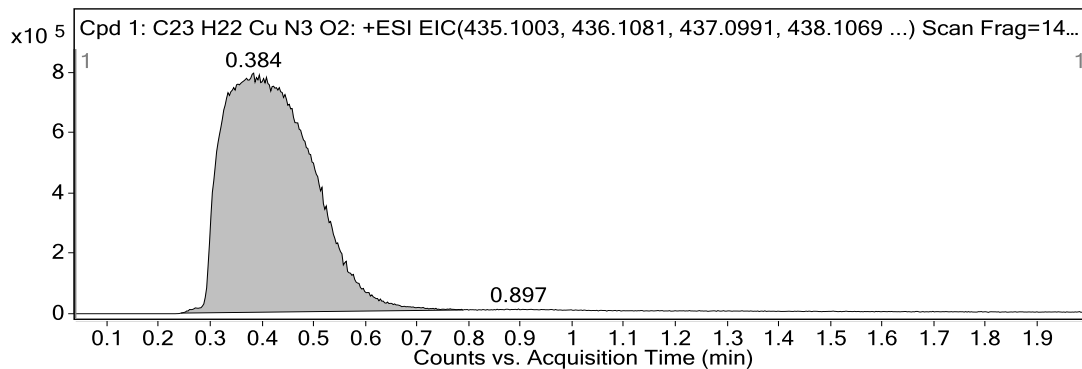
<b>Data File</b>	MJ_MS_2DMSO_MK_140V_pos1.d	<b>Sample Name</b>	2DMSO
<b>Sample Type</b>	Sample	<b>Position</b>	Vial 6
<b>Instrument Name</b>	DE1517B001	<b>User Name</b>	
<b>Acq Method</b>	Odredjivanje MM_MK_140V_pos.m	<b>Acquired Time</b>	11/29/2019 2:08:25 PM
<b>IRM Calibration Status</b>	Success	<b>DA Method</b>	EURL_target_PEST017_NNP_10V.m
<b>Comment</b>			

<b>Sample Group</b>		<b>Info.</b>	
<b>Stream Name</b>	LC 1	<b>Acquisition SW Version</b>	6200 series TOF/6500 series Q-TOF B.06.01 (B6157)

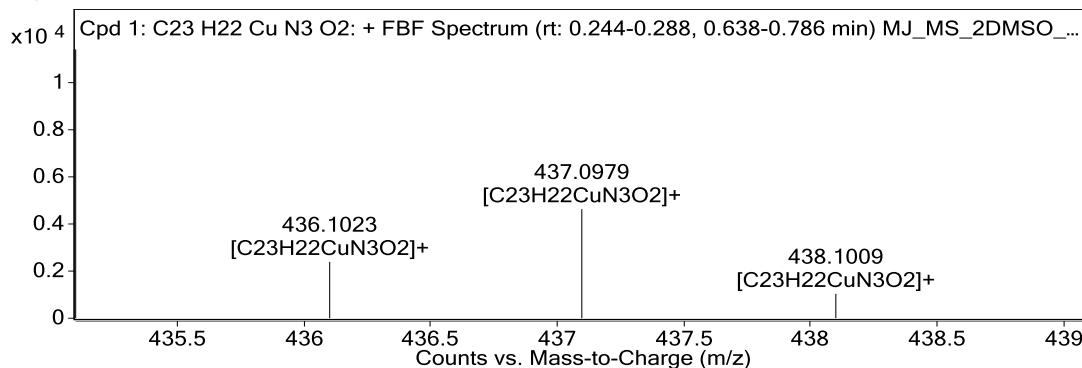
## Compound Table

Compound Label	RT	Mass	Abund	Formula	Tgt Mass	Diff (ppm)
Cpd 1: C23 H22 Cu N3 O2	0.384	435.0998	11398	C23 H22 Cu N3 O2	435.1008	-2.37

Compound Label	m/z	RT	Algorithm	Mass
Cpd 1: C23 H22 Cu N3 O2	435.0993	0.384	Find By Formula	435.0998

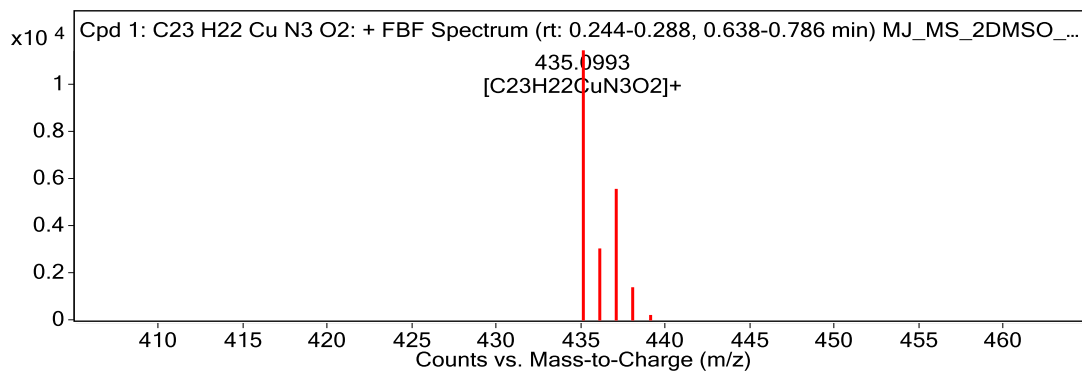


## MS Spectrum



## MS Zoomed Spectrum

# Qualitative Compound Report



## MS Spectrum Peak List

m/z	z	Abund	Formula	Ion
435.0993	1	11398.25	C <sub>23</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
436.1023	1	2389.83	C <sub>23</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
437.0979	1	4626.48	C <sub>23</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
438.1009	1	1034.49	C <sub>23</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>
439.1027	1	31.45	C <sub>23</sub> H <sub>22</sub> CuN <sub>3</sub> O <sub>2</sub>	M <sup>+</sup>

--- End Of Report ---